LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

Signal : TIC: BP013312.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	21	24	33	rVB	117538	160591	1.54%	0.115%
2	3.769	94	99	110	rBV2	180649	301610	2.89%	0.217%
3	3.863	111	115	122	rVV	154923	207463	1.99%	0.149%
4	3.987	131	136	142	rVB	115526	181606	1.74%	0.131%
5	4.087	149	153	158	rVB	51818	68596	0.66%	0.049%
6	4.463	213	217	224	rVB4	28250	41932	0.40%	0.030%
7	4.652	245	249	253	rVV	64685	90114	0.86%	0.065%
8	4.699	253	257	269	rVB	154688	226558	2.17%	0.163%
9	5.016	305	311	323	rVV2	391260	759408	7.27%	0.546%
10	5.116	325	328	337	rVB2	23895	39389	0.38%	0.028%
11	6.540	565	570	575	rBV2	38165	63330	0.61%	0.046%
12	6.610	576	582	587	rVB	78782	126401	1.21%	0.091%
13	7.098	658	665	680	rVB	1780836	2917313	27.92%	2.097%
14	7.263	686	693	701	rBV	1504836	2335821	22.35%	1.679%
15	7.469	721	728	736	rVB	1924293	2961563	28.34%	2.129%
16	7.551	736	742	746	rBV2	25453	43900	0.42%	0.032%
17	7.940	801	808	815	rBV	2496261	3897195	37.30%	2.802%
18	8.651	922	929	940	rBV	1695353	2711435	25.95%	1.949%
19	8.769	943	949	958	rVB	164163	265322	2.54%	0.191%
20	9.104	999	1006	1015	rBV	1649107	2768783	26.50%	1.990%
21	9.263	1029	1033	1040	rVB7	15753	30190	0.29%	0.022%
22	9.504	1067	1074	1088	rVB2	53097	101884	0.98%	0.073%
23	9.834	1121	1130	1144	rBV	1391950	2311476	22.12%	1.662%
24	9.981	1151	1155	1163	rVB2	30436	55496	0.53%	0.040%
25	10.187	1186	1190	1196	rVB4	16417	30657	0.29%	0.022%
26	10.375	1214	1222	1236	rBV	2021220	3387198	32.42%	2.435%
27	10.751	1277	1286	1293	rBV	3133791	5180972	49.58%	3.725%
28	10.798	1293	1294	1303	rVB	28461	40530	0.39%	0.029%
29	10.892	1303	1310	1327	rBV	1009502	1856362	17.77%	1.335%
30	11.287	1370	1377	1384	rBV	17537	44901	0.43%	0.032%
31	11.545	1415	1421	1430	rBV6	15485	34729	0.33%	0.025%
32	12.334	1550	1555	1561	rVB	35286	56439	0.54%	0.041%
33	13.104	1682	1686	1691	rVB3	22598	33054	0.32%	0.024%
34	13.398	1727	1736	1740	rBV2	41517	70584	0.68%	0.051%
35	13.904	1817	1822	1828	rVB4	45910	69193	0.66%	0.050%
36	13.986	1828	1836	1845	rBV	2859631	4162974	39.84%	2.993%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

37	14.110	1851	1857	1863	rVB8	24764	38476	0.37%	0.028%
38	14.169	1863	1867	1873	rBV	42131	60039	0.57%	0.043%
39	14.280	1879	1886	1898	rVV	3304090	4998978	47.84%	3.594%
40	14.469	1914	1918	1923	rVB	30537	41522	0.40%	0.030%
41	14.580	1931	1937	1944	rVV	3684862	5611444	53.70%	4.034%
42	14.645	1944	1948	1959	rVB	85901	146312	1.40%	0.105%
43	14.786	1965	1972	1990	rBV	925023	1635691	15.65%	1.176%
44	14.939	1994	1998	2002	rVV4	60298	105392	1.01%	0.076%
45	14.986	2002	2006	2010	rVV3	86187	163368	1.56%	0.117%
46	15.033	2010	2014	2021	rVB2	50931	92086	0.88%	0.066%
47	15.245	2045	2050	2056	rVB9	19507	33186	0.32%	0.024%
48	15.369	2066	2071	2074	rBV6	14717	28259	0.27%	0.020%
49	15.580	2099	2107	2112	rBV	4137482	5975183	57.18%	4.296%
50	15.633	2112	2116	2121	rVB	133720	196599	1.88%	0.141%
51	15.698	2121	2127	2135	rBV	707486	985011	9.43%	0.708%
52	15.945	2160	2169	2179	rVV	2389352	3236945	30.98%	2.327%
53	16.080	2188	2192	2198	rVB4	21926	35176	0.34%	0.025%
54	16.151	2198	2204	2210	rVB4	14250	31961	0.31%	0.023%
55	16.427	2244	2251	2254	rBV5	15115	34683	0.33%	0.025%
56	16.474	2254	2259	2262	rVV	97560	141175	1.35%	0.101%
57	16.510	2262	2265	2271	rVB	112666	149334	1.43%	0.107%
58	16.722	2298	2301	2310	rBV9	31337	51395	0.49%	0.037%
59	17.039	2352	2355	2359	rVB2	29610	33963	0.33%	0.024%
60	17.092	2359	2364	2367	rBV6	17940	31386	0.30%	0.023%
61	17.157	2371	2375	2380	rVB	56349	83336	0.80%	0.060%
62	17.227	2383	2387	2394	rVB8	24813	42234	0.40%	0.030%
63	17.345	2400	2407	2411	rBV	4212268	6147716	58.84%	4.420%
64	17.386	2411	2414	2418	rVV	991604	1331575	12.74%	0.957%
65	17.439	2418	2423	2436	rVV2	4183205	6316021	60.45%	4.541%
66	17.533	2436	2439	2445	rVB3	41114	66763	0.64%	0.048%
67	17.616	2450	2453	2459	rVB2	27803	41746	0.40%	0.030%
68	17.745	2468	2475	2480	rBV	119390	187584	1.80%	0.135%
69	17.998	2515	2518	2523	rVB5	47910	56705	0.54%	0.041%
70	18.098	2532	2535	2540	rVB5	29902	43030	0.41%	0.031%
71	18.157	2541	2545	2548	rBV2	40648	54673	0.52%	0.039%
72	18.216	2549	2555	2559	rBV2	744247	1065424	10.20%	0.766%
73	18.404	2581	2587	2594	rBV4	198964	339511	3.25%	0.244%
74	18.463	2594	2597	2600	rVB2	56921	59045	0.57%	0.042%
75	18.545	2607	2611	2616	rBV3	107888	164772	1.58%	0.118%
76	18.621	2621	2624	2628	rVV	146545	174216	1.67%	0.125%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Title : SVOA CALIBRATION

77	18.686	2632	2635	2638	rVV2	77810	106064	1.02%	0.076%
78	18.721	2639	2641	2646	rVV5	62772	98989	0.95%	0.071%
79	18.780	2647	2651	2659	rVB	255402	321330	3.08%	0.231%
80	19.116	2706	2708	2711	rVB2	85656	88057	0.84%	0.063%
81	19.351	2738	2748	2755	rBV	1964065	2771063	26.52%	1.992%
82	19.445	2760	2764	2769	rBV2	249266	348275	3.33%	0.250%
83	19.586	2786	2788	2796	rVB2	94018	109225	1.05%	0.079%
84	19.674	2797	2803	2806	rBV	6908133	9263530	88.66%	6.659%
85	19.704	2806	2808	2815	rVB	1827743	1954897	18.71%	1.405%
86	19.821	2824	2828	2834	rBV	1134001	1378813	13.20%	0.991%
87	20.015	2857	2861	2862	rBV	77393	114074	1.09%	0.082%
88	20.180	2884	2889	2895	rVB2	444355	796777	7.63%	0.573%
89	20.280	2902	2906	2910	rBV	341137	395831	3.79%	0.285%
90	21.080	3039	3042	3045	rBV	232243	287283	2.75%	0.207%
91	21.127	3045	3050	3054	rBV2	1511263	2535537	24.27%	1.823%
92	21.433	3095	3102	3106	rBV2	6254938	10448875	100.00%	7.512%
93	21.468	3106	3108	3115	rVB	1302755	1676585	16.05%	1.205%
94	22.068	3207	3210	3220	rVB3	686338	1214452	11.62%	0.873%
95	23.151	3386	3394	3399	rBV2	1276337	3595718	34.41%	2.585%
96	23.692	3482	3486	3490	rBV	643903	1143057	10.94%	0.822%
97	23.751	3490	3496	3502	rBV2	4328091	8911449	85.29%	6.406%
98	23.915	3517	3524	3530	rBV2	3881193	8018102	76.74%	5.764%
99	24.527	3623	3628	3638	rVB	1298415	2779562	26.60%	1.998%
100	26.439	3947	3953	3959	rBV	1215576	3078384	29.46%	2.213%

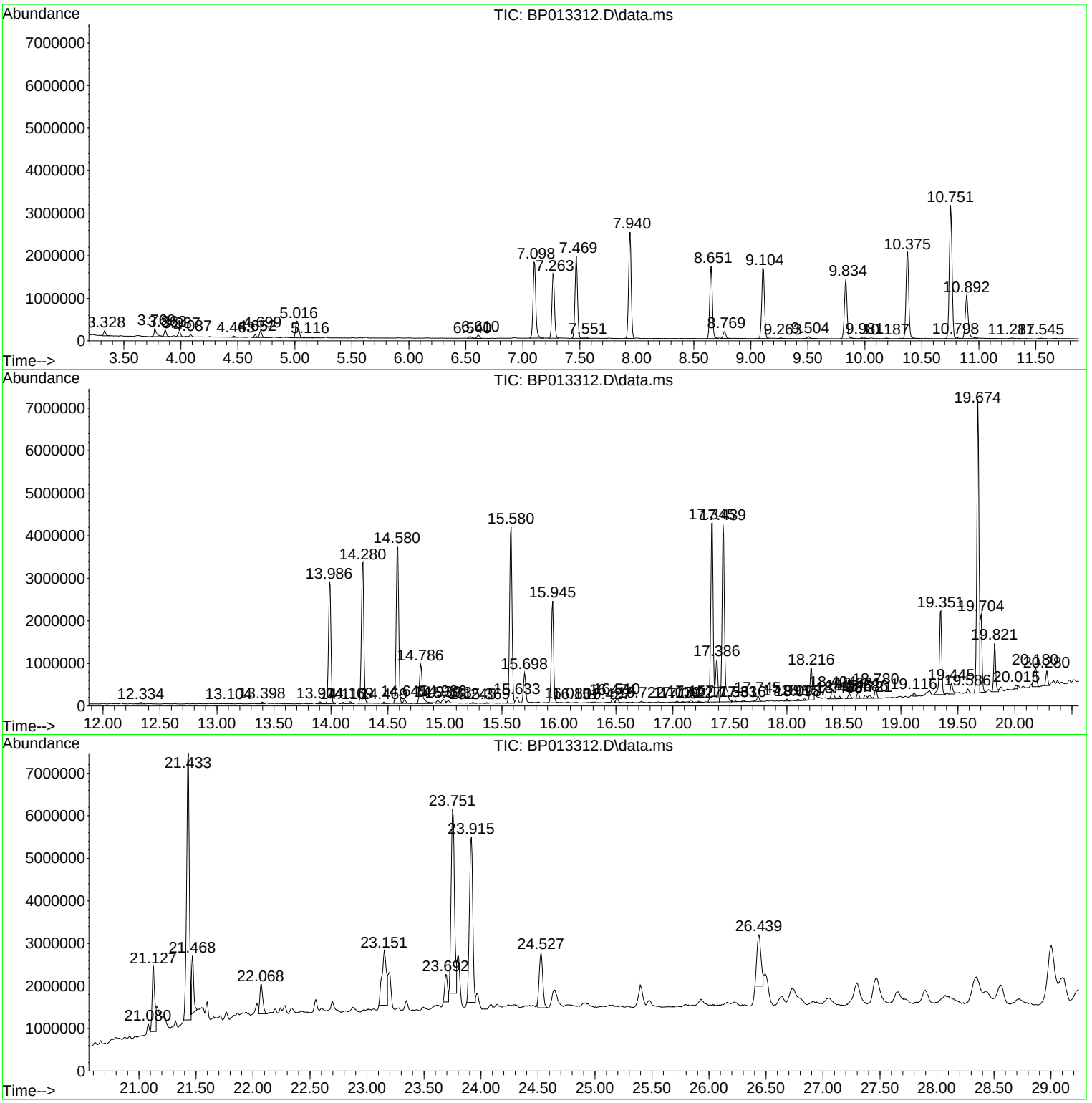
Sum of corrected areas: 139102813

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

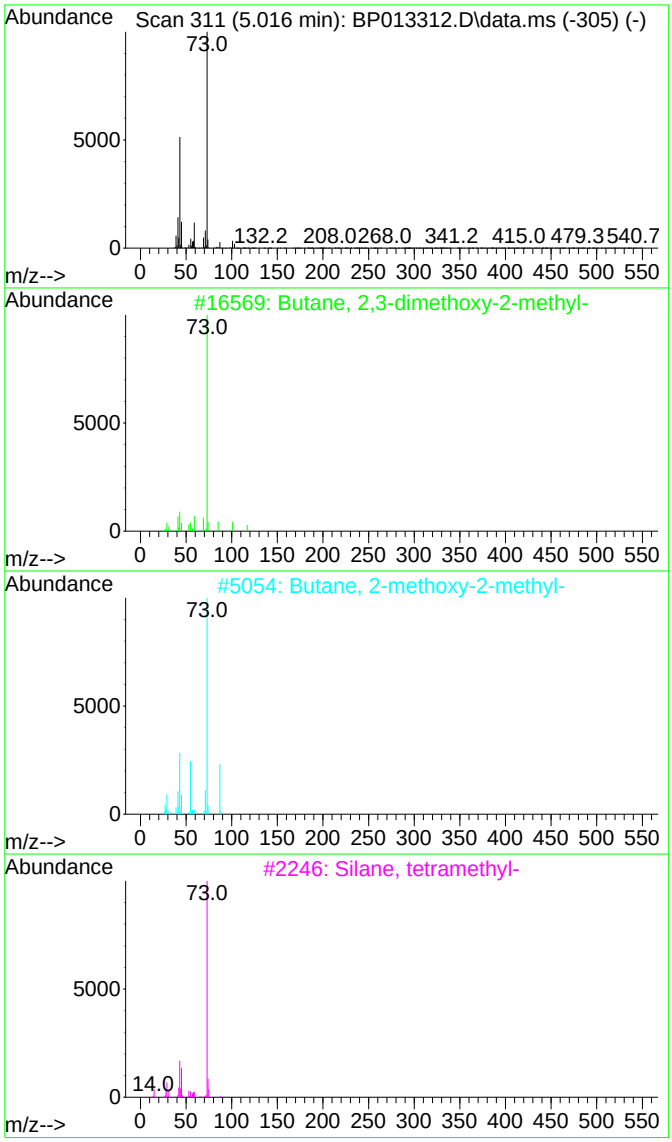
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

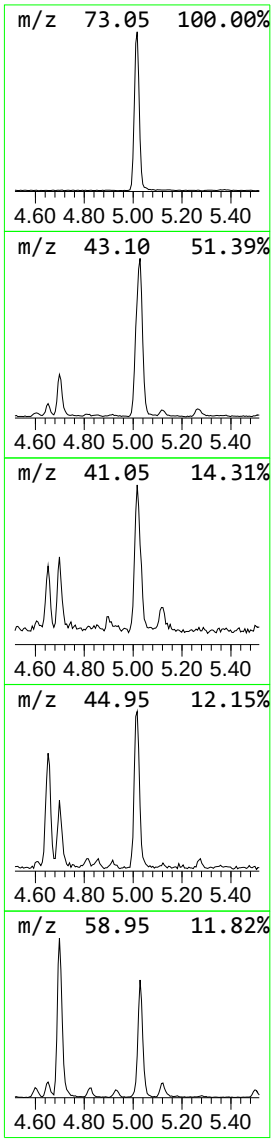
****

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	3.90 ng/ul	759408	1,4-Dichlorobenzene-d4	7.940

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethoxy-2-methyl-	132	C7H16O2	074421-00-4	45
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	33
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	32
4	2-Pentene, 2-methoxy-	100	C6H12O	061142-47-0	23
5	4-Ethoxy-2-butanone	116	C6H12O2	060044-74-8	17





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

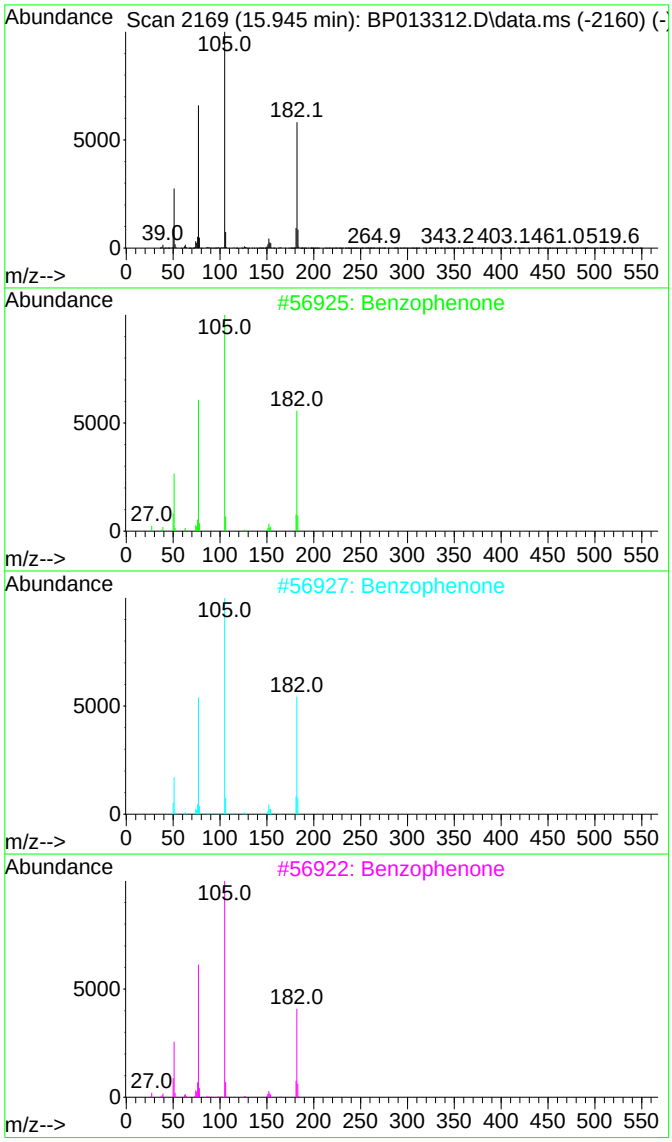
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

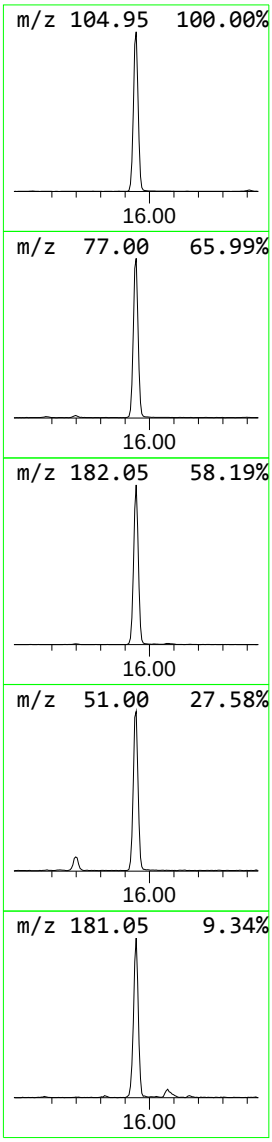
****

Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	11.54 ng/ul	3236950	Acenaphthene-d10	14.586

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	95
3	Benzophenone	182	C13H10O	000119-61-9	94
4	Benzophenone	182	C13H10O	000119-61-9	91
5	Benzophenone	182	C13H10O	000119-61-9	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

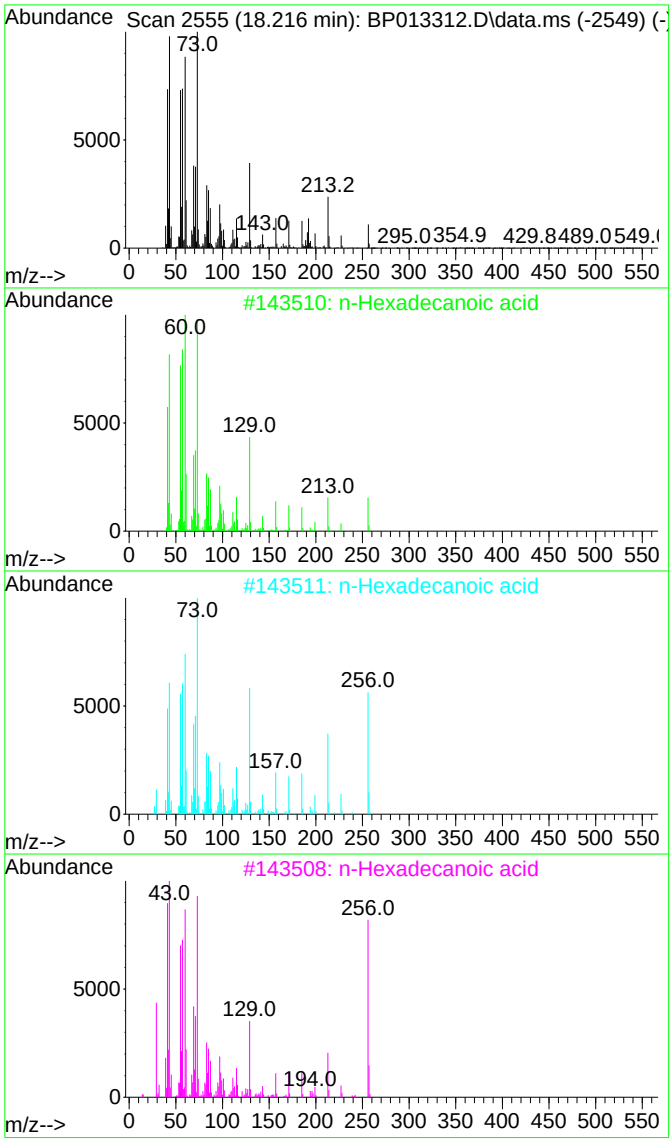
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

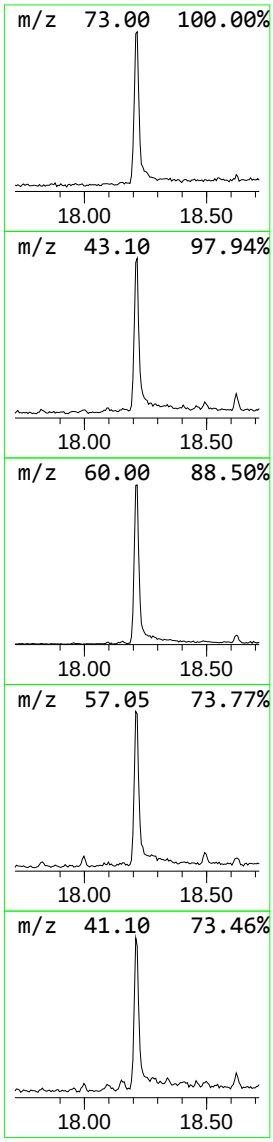
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	3.47 ng/ul	1065420	Phenanthrene-d10	17.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

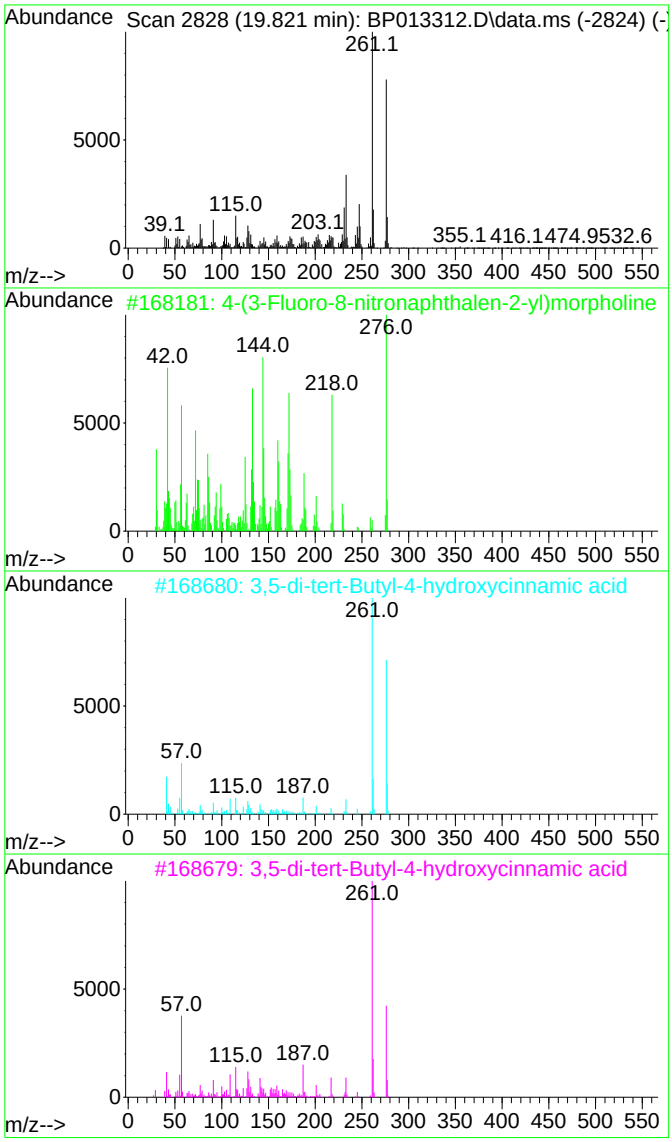
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

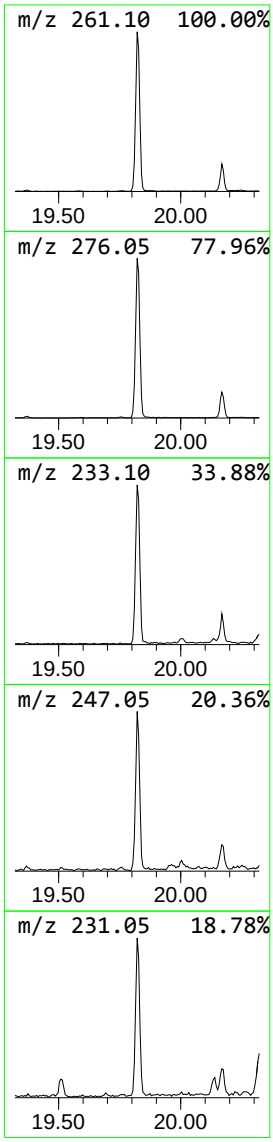
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 4-(3-Fluoro-8-nitronaphthal... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.821	2.64 ng/ul	1378810	Chrysene-d12	21.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(3-Fluoro-8-nitronaphthalen-2-...	276	C14H13FN2O3	1000409-37-1	92
2	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	68
3	3,5-di-tert-Butyl-4-hydroxycinna...	276	C17H24O3	022014-01-3	68
4	(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	68
5	5H-Furo[3,2-g][1]benzopyran-5-on...	276	C14H12O6	000668-10-0	62





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

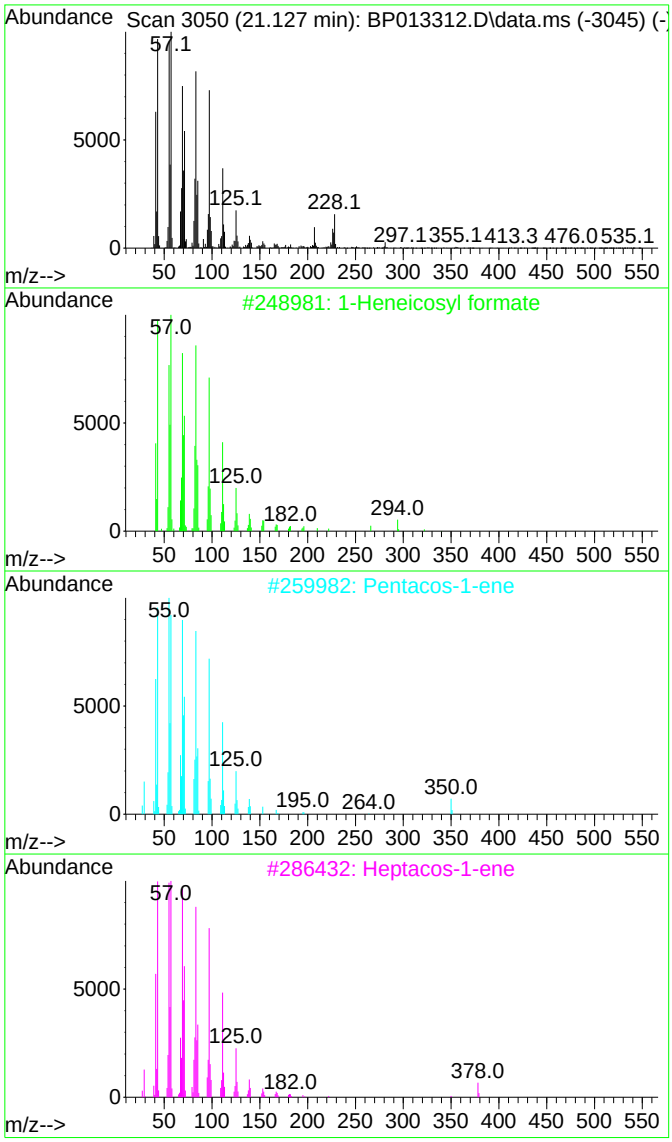
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

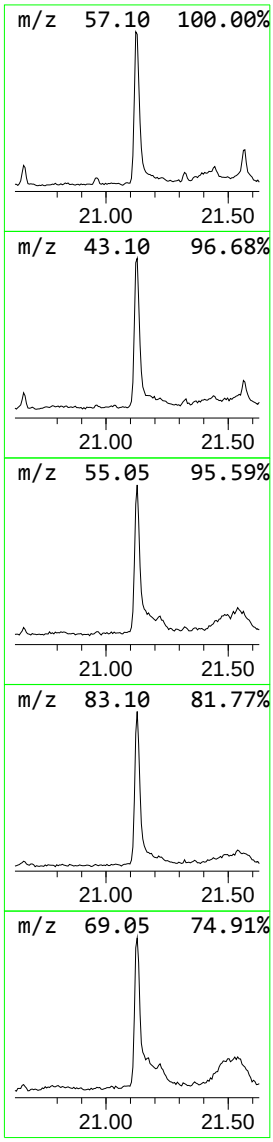
****

Peak Number 5 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	4.85 ng/ul	2535540	Chrysene-d12	21.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2	Pentacos-1-ene	350	C25H50	016980-85-1	94
3	Heptacos-1-ene	378	C27H54	015306-27-1	94
4	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

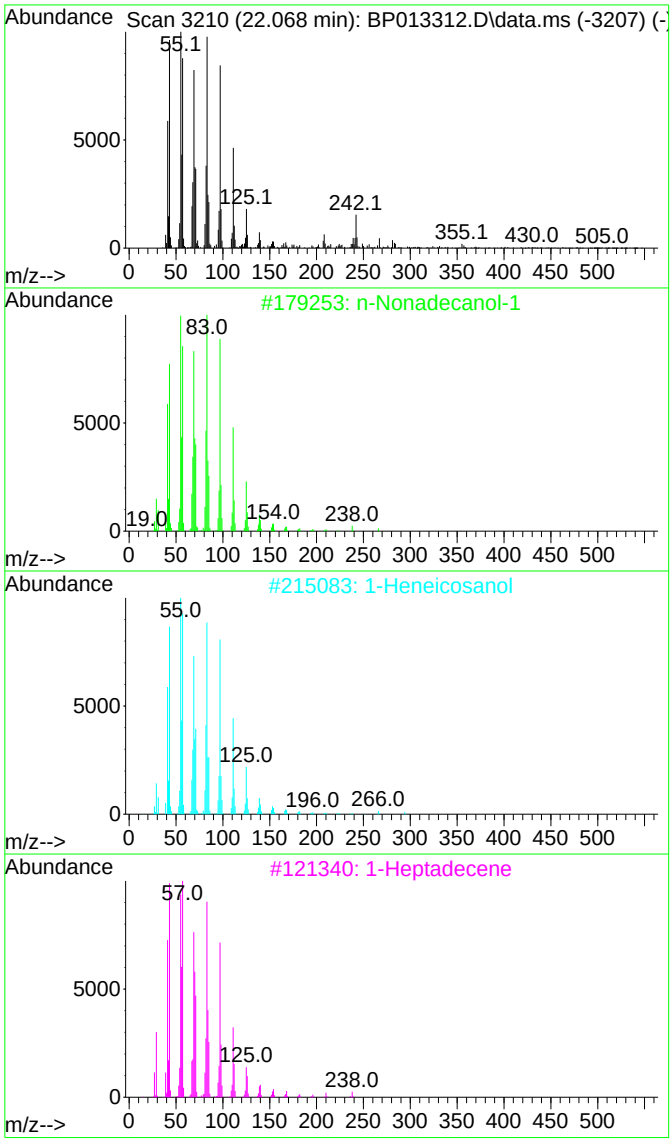
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

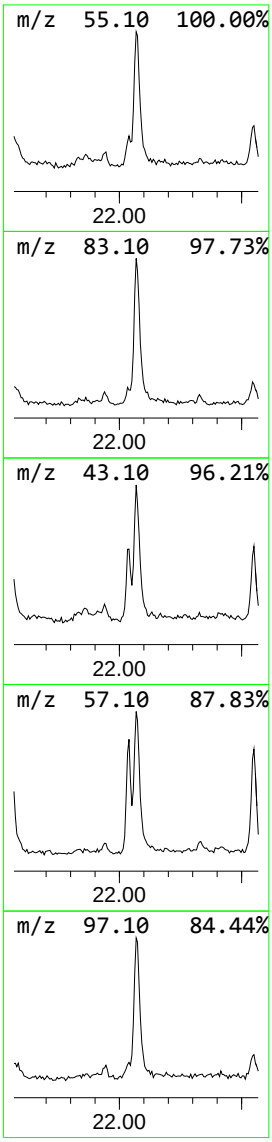
****

Peak Number 6 n-Nonadecanol-1 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.068	2.32 ng/ul	1214450	Chrysene-d12	21.433

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	95
3	1-Heptadecene	238	C17H34	006765-39-5	95
4	Behenic alcohol	326	C22H46O	000661-19-8	94
5	n-Heptadecanol-1	256	C17H36O	001454-85-9	93





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Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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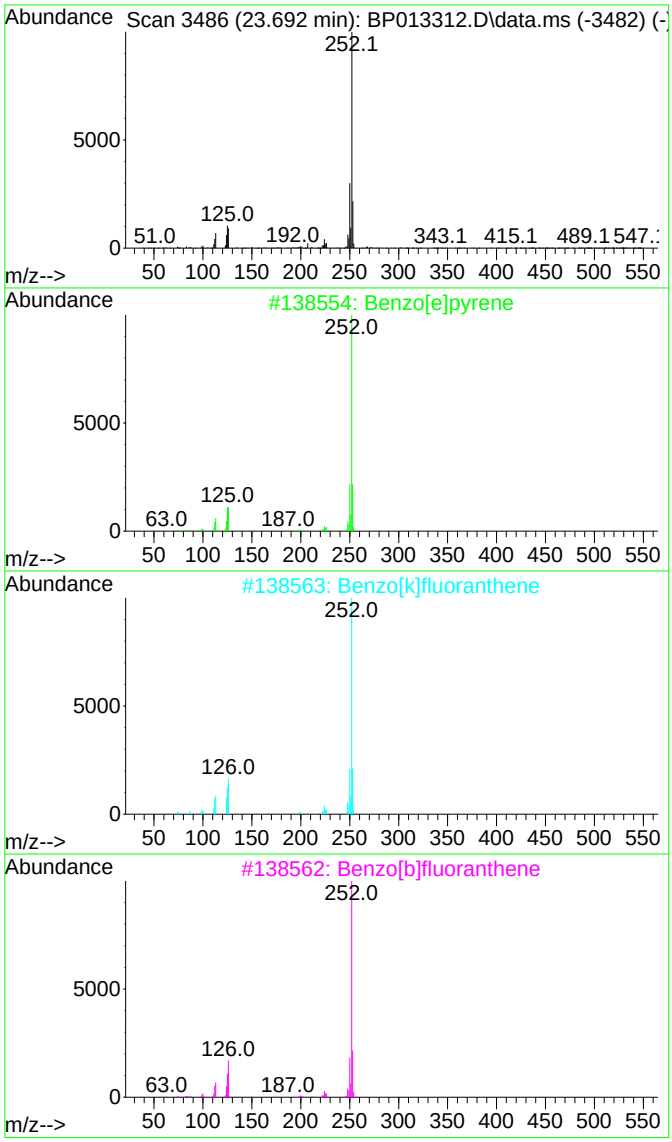
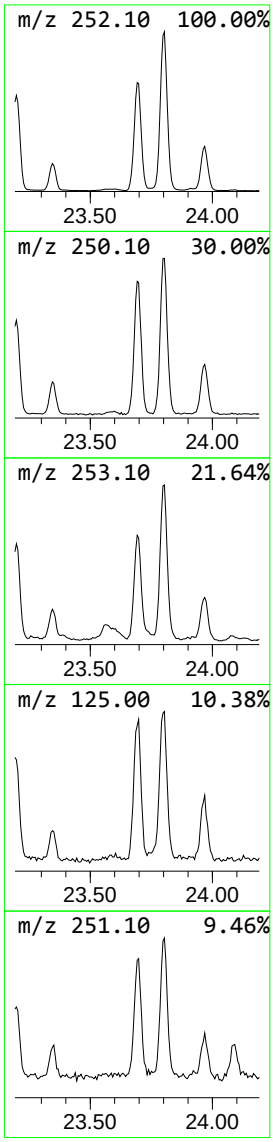
Peak Number 7 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.692	2.85 ng/ul	1143060	Perylene-d12	23.915

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
5	Benzo[a]pyrene	252	C20H12	000050-32-8	93





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Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

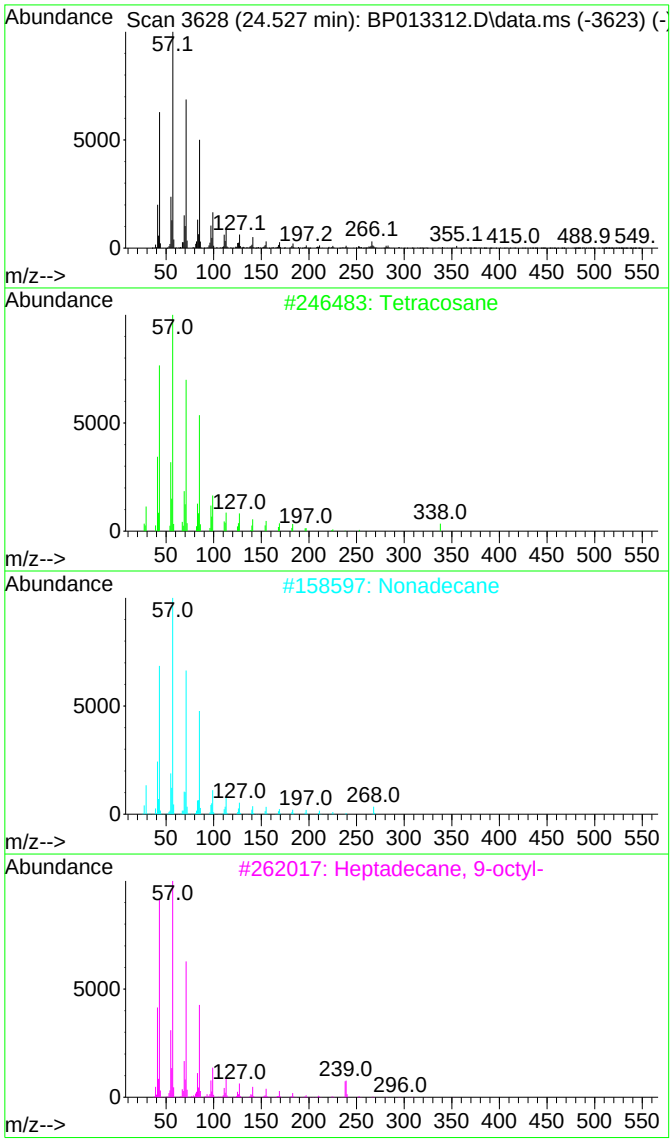
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

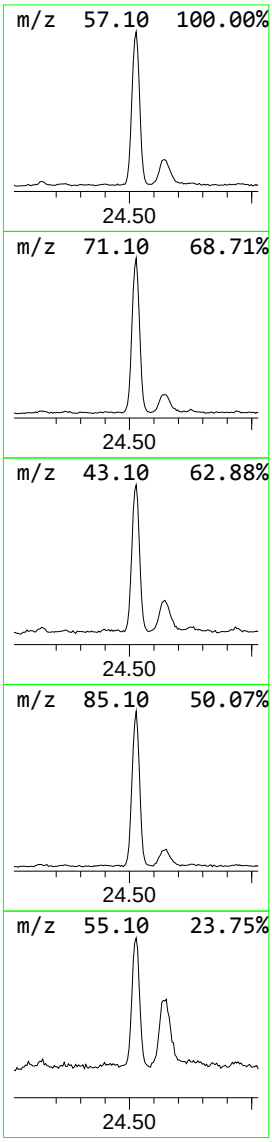
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	6.93 ng/ul	2779560	Perylene-d12	23.915

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	94
2	Nonadecane	268	C19H40	000629-92-5	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
4	Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5	Tetracosane	338	C24H50	000646-31-1	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

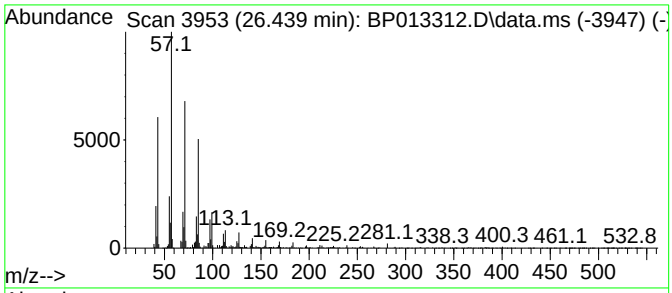
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

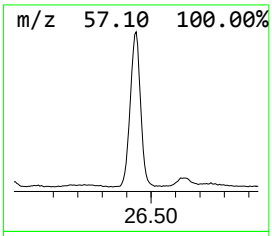
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

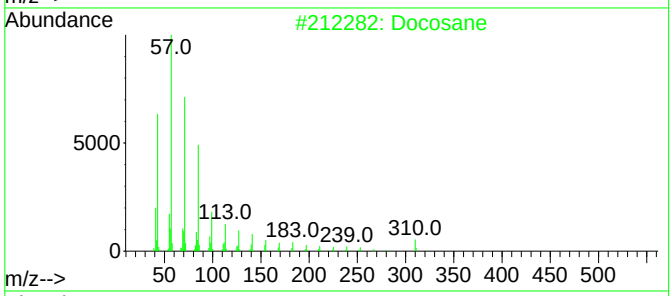
****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

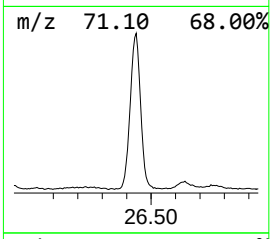
R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	7.68 ng/ul	3078380	Perylene-d12	23.915

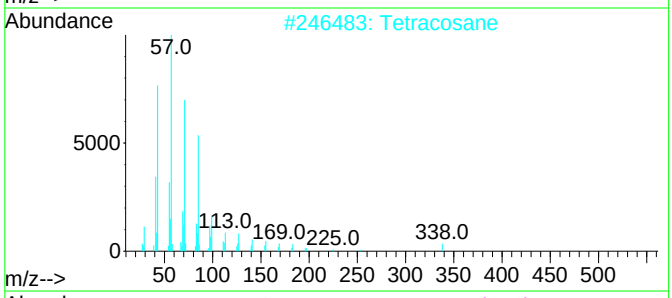
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane	310	C22H46	000629-97-0	97
2	Tetracosane	338	C24H50	000646-31-1	95
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	94
5	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94

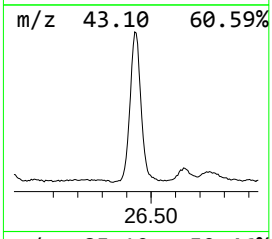


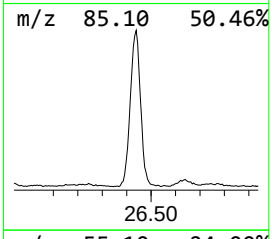


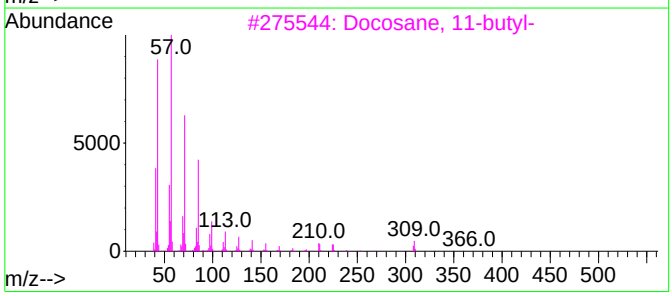


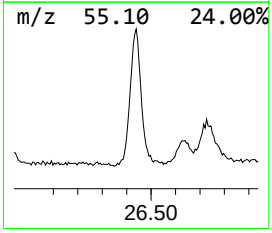












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013312.D
Acq On : 24 Dec 2022 22:35
Operator : CG/JU
Sample : N6016-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	5.016	3.9	ng/ul	759408	1	7.940	3897200	20.0
Benzophenone	15.945	11.5	ng/ul	3236950	3	14.586	5611440	20.0
n-Hexadecanoic ...	18.216	3.5	ng/ul	1065420	4	17.345	6147720	20.0
4-(3-Fluoro-8-n...	19.821	2.6	ng/ul	1378810	5	21.433	10448900	20.0
1-Heneicosyl fo...	21.127	4.8	ng/ul	2535540	5	21.433	10448900	20.0
n-Nonadecanol-1	22.068	2.3	ng/ul	1214450	5	21.433	10448900	20.0
Benzo[e]pyrene	23.692	2.9	ng/ul	1143060	6	23.915	8018100	20.0
(DEL) Alkane: S...	24.527	6.9	ng/ul	2779560	6	23.915	8018100	20.0
(DEL) Alkane: S...	26.439	7.7	ng/ul	3078380	6	23.915	8018100	20.0